## **STN Columbus**

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NEWS
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NEWS
                 CA/CAplus pre-1967 chemical substance index entries enhanced
NEWS
         DEC 18
                 with preparation role
                 CA/CAplus patent kind codes updated
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         DEC 18
                 MARPAT to CA/CAplus accession number crossover limit increased
         DEC 18
NEWS
                 to 50,000
                 MEDLINE updated in preparation for 2007 reload
NEWS
         DEC 18
         DEC 27
                 CA/CAplus enhanced with more pre-1907 records
NEWS
      7
                 CHEMLIST enhanced with New Zealand Inventory of Chemicals
NEWS
         JAN 08
                 CA/CAplus Company Name Thesaurus enhanced and reloaded
         JAN 16
NEWS
     9
NEWS 10
         JAN 16
                 IPC version 2007.01 thesaurus available on STN
                 WPIDS/WPINDEX/WPIX enhanced with IPC 8 reclassification data
NEWS 11
         JAN 16
                 CA/CAplus updated with revised CAS roles
NEWS 12
         JAN 22
                 CA/CAplus enhanced with patent applications from India
NEWS 13
         JAN 22
                 PHAR reloaded with new search and display fields
         JAN 29
NEWS 14
                 CAS Registry Number crossover limit increased to 300,000 in
NEWS 15
         JAN 29
                 multiple databases
                 PATDPASPC enhanced with Drug Approval numbers
NEWS 16
         FEB 15
                 RUSSIAPAT enhanced with pre-1994 records
NEWS 17
         FEB 15
                 KOREAPAT enhanced with IPC 8 features and functionality
NEWS 18
         FEB 23
                 MEDLINE reloaded with enhancements
NEWS 19
         FEB 26
                 EMBASE enhanced with Clinical Trial Number field
NEWS 20
         FEB 26
                 TOXCENTER enhanced with reloaded MEDLINE
NEWS 21
         FEB 26
NEWS 22
         FEB 26
                 IFICDB/IFIPAT/IFIUDB reloaded with enhancements
                 CAS Registry Number crossover limit increased from 10,000
         FEB 26
NEWS 23
                  to 300,000 in multiple databases
                 WPIDS/WPIX enhanced with new FRAGHITSTR display format
NEWS 24
         MAR 15
NEWS 25
                 CASREACT coverage extended
         MAR 16
NEWS 26
                 MARPAT now updated daily
         MAR 20
         MAR 22
                 LWPI reloaded
NEWS 27
NEWS 28
         MAR 30
                 RDISCLOSURE reloaded with enhancements
                 INPADOCDB will replace INPADOC on STN
NEWS 29
         MAR 30
              NOVEMBER 10 CURRENT WINDOWS VERSION IS V8.01c, CURRENT
NEWS EXPRESS
              MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP)
              AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.
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                                                 SINCE FILE
                                                                 TOTAL
                                                               SESSION
                                                      ENTRY
                                                       0.21
                                                                  0.21
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## http://www.cas.org/ONLINE/UG/reqprops.html

```
=> e fludarabine/cn
                      FLUDALANINE/CN
               1
                      FLUDARA/CN
E2
               1
                 --> FLUDARABINE/CN
E3
                      FLUDARABINE 5'-MONOPHOSPHATE/CN
E4
               1
                      FLUDARABINE PHOSPHATE/CN
E5
               1
                      FLUDARABINE TRIPHOSPHATE/CN
E6
               1
                      FLUDARENE/CN
E7
               1
E8
                      FLUDAZONIUM CHLORIDE/CN
                      FLUDELONE/CN
E9
               1
E10
               1
                      FLUDEMIL/CN
                      FLUDENT/CN
E11
               1
E12
               1
                      FLUDEOXYGLUCOSE (18F)/CN
=> s e3
               1 FLUDARABINE/CN
L1
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2007 ACS on STN 21679-14-1 REGISTRY
L1
RN
     Entered STN: 16 Nov 1984
ED
      9H-Purin-6-amine, 9-\beta-D-arabinofuranosyl-2-fluoro- (9CI)
                                                                          (CA INDEX
CN
OTHER CA INDEX NAMES:
     Adenine, 9-\beta-D-arabinofuranosyl-2-fluoro- (8CI)
OTHER NAMES:
     2-Fluoro Ara-A
CN
     2-Fluoro-9-β-D-arabinofuranosyladenine
     2-Fluoroadenine arabinoside
CN
      9-\beta-D-Arabinofuranosyl-2-fluoroadenine
CN
      9-β-D-Arabinosyl-2-fluoroadenine
CN
CN
     F-ara-A
CN
     Fludarabine
     NSC 118218
NSC 118218H
CN
CN
      STEREOSEARCH
FS
      C10 H12 F N5 O4
MF
CI
      COM
        N Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DRUGU, EMBASE, HSDB*, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE,
      STN Files:
        MRCK*, PHAR, PROMT, PS, RTECS*, SYNTHLINE, TOXCENTER, USAN, USPAT2,
        USPATFULL
           (*File contains numerically searchable property data)
                         EINECS**, WHO
           (**Enter CHEMLIST File for up-to-date regulatory information)
```

Absolute stereochemistry.

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1412 REFERENCES IN FILE CA (1907 TO DATE)
38 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1418 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file merck
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 7.35 7.56

FULL ESTIMATED COST

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FILE 'REGISTRY' ENTERED AT 17:19:02 ON 30 MAR 2007 E FLUDARABINE/CN

L1

1 S E3

FILE 'MRCK' ENTERED AT 17:19:45 ON 30 MAR 2007

=> s 11

L2 1 L1

=> d all

L2 ANSWER 1 OF 1 MRCK COPYRIGHT (C) 2007 Merck and Co., Inc., Whitehouse Station, New Jersey, USA. All rights reserved. on STN

MERCK Number (MNO): 4152

CAS Registry No. (RN): 21679-14-1 MERCK Index Name (MIN): Fludarabine

CA Index Name (CN):  $9-\beta-D$ -Arabinofuranosyl-2-fluoro-9H-purin-6-amine

Synonym(s) (CN):  $9-\beta-D$ -arabinofuranosyl-2-fluoroadenine; 2-fluorovidarabine; 2-fluoro-9- $\beta$ -D-

arabinofuranosyladenine; 2-F-araA Drug Code(s) (CN): NSC-118218; NSC-118218-H

Molecular Form. (MF): C10 H12 F N5 O4

Wgt Composition (COMP): C 42.11%, H 4.24%, F 6.66%, N 24.55%, O 22.44%.

Molecular Weight (MW): 285.23

References (RE): Adenosine deaminase-resistant purine nucleoside antimetabolite. Prepn and in vitro cytotoxicity: J. A. Montgomery, K. Hewson, J. Med. Chem. 12, 498 (1969). Improved prepn: J. A. Montgomery et al., J. Heterocycl. Chem. 16, 157 (1979); J. A. Montgomery, US 4210745 (1980 to U.S. Dept. Health, Education and Welfare). Inhibition of DNA synthesis and in vivo antileukemic activity: R. W. Brockman et al., Biochem. Pharmacol. 26, 2193 (1977). Metabolized to 5'-monophosphate: R. W. Brockman et al., Cancer Res. 40, 3610 (1980). HPLC determn in human leukemia cells: V. Gandhi et al., J. Chromatogr. 413, 293 (1987). Prepn

of 5'-monophosphate: J. A. Montgomery, A. T. Shortnacy, US 4357324 (1982 to U.S. Dept. of Health and Human Services). Pharmacokinetics in humans: M. R. Hersh et al., Cancer Chemother. Pharmacol. 17, 277 (1986). Evaluation of therapeutic efficacy and CNS toxicity in acute refractory leukemia: R. P. Warrell, Jr., E. Berman, J. Clin. Oncol. 4, 74 (1986); H. G. Chun et al., Cancer Treat. Rep. 70, 1225 (1986). Series of articles on pharmacology and therapeutic use: Semin. Oncol. 17, Suppl. 8, 1-78 (1990).

Absolute stereochemistry.

Melting Point (MP):

Value MP deg C ==== 260

Optical Rotatory Power (ORP):

| Value  | Temp. | Spectral | Note                 |
|--------|-------|----------|----------------------|
| ORP    | ORP.T | Line     |                      |
| deg    | deg C | ORP.SL   |                      |
| 17 2.5 | 25    | D        | (c = 0.1 in ethanol) |

UV Spectrum (UVS):

Other Properties (OCPP): .

Crystals from ethanol + water, mp 260°. [ $\alpha$ ]D25 +17  $\pm 2.5^{\circ}$  (c = 0.1 in ethanol) . uv max (pH 1, pH 7, pH 13): 262 , 261 , 262 nm ( $\epsilon$  x 10-3 13.2, 14.8, 15.0) . Sparingly sol in water, organic solvents.

```
== DERIVATIVE == (1): 5'-Monophosphate

CAS Registry No. (RN.DRV): 75607-67-9

Synonym(s) (CN.DRV): 2-F-ara-AMP

Drug Code(s) (CN.DRV): NSC-328002; NSC-312887

Trade Name(s) (CN.DRV): Fludara (Schering AG)

Molecular Form. (MF.DRV): C10 H13 F N5 O7 P

Wgt Composition (COMP.DRV): C 32.89%, H 3.59%, F 5.20%, N 19.18%, O 30.67%, P

8.48%.

Molecular Weight (MW.DRV): 365.21
```

## Absolute stereochemistry.

Other Properties (OCPP.DRV): Sol in water.

Therapeutic Codes (THER):

Phosphate as antineoplastic.
Referenced Patent (RPN):

US4210745; US4357324

=> file drugu
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 4.61 12.17

FULL ESTIMATED COST

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>>> FILE COVERS 1983 TO DATE <>>
>>> THESAURUS AVAILABLE IN /CT <>>

=> s l1

L3 1541 L1

=> d 1541

L3 ANSWER 1541 OF 1541 DRUGU COPYRIGHT 2007 THE THOMSON CORP on STN

AN 9572 DRUGU FS Registry DDRN FLUDARABI DDN FLUDARABINE RN 21679-14-1

CYTOSTATICS

SS AMIDINE, CYCLIC; HH-LINKED-CX; PURINE; NUCLEOSIDE; FLUORINE

=> log y

CT

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 8.42 20.59

STN INTERNATIONAL LOGOFF AT 17:28:18 ON 30 MAR 2007